

PIRLS: Poisson Iteratively Reweighted Least Squares Computer Program for Additive, Multiplicative, Power, and Non-linear Models

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Abstract

A computer program to estimate Poisson regression coefficients, standard errors, Pearson χ^2 and deviance goodness of fit (and residuals), leverages, and Freeman-Tukey, standardized, and deletion residuals for additive, multiplicative, power, and non-linear models is described. Data used by the program must be stored and input from a disk file. The output file contains an optional list of the input data, observed and fitted count data (deaths or cases), coefficients, standard errors, relative risks and 95% confidence intervals, variance-covariance matrix, correlation matrix, goodness of fit statistics, and residuals for regression diagnostics.

Key Words: Poisson Regression, Iteratively Reweighted Least Squares, Regression Diagnostics, Matrix Manipulation.

1 INTRODUCTION

1.1 Poisson Distribution

Poisson regression models are commonly used for modeling risk factors and other covariates when the rates of disease are small, i.e., cancer [1, 2, 3, 4, 5]. An area of research in which Poisson regression of cancer rates is commonly employed is the investigation of radiation-induced cancer among Hiroshima and Nagasaki atomic bomb survivors [6, 7, 8] and individuals occupationally [9], medically [10] and environmentally [11] exposed to ionizing radiation. Poisson regression has also been used to investigate the effects of information bias on lifetime risks [12], diagnostic misclassification [13, 14], and dose-response curves for radiation-induced chromosome aberrations [15].

Several computer programs have been developed that perform Poisson regression [16, 17, 18, 19, 20]. While the PREG program [3] can fit additive, multiplicative, and non-linear models, it does not provide the capability to fit power models nor perform regression diagnostics. This report documents the PIRLS Poisson regression algorithm for fitting additive, multiplicative, power, and non-linear models and regression diagnostics using leverages of the Hat matrix, deletion and standardized residuals.

The critical data required to apply the computer code are:

- Number of cases for each stratum
- Number of person-years for each stratum
- Design matrix of independent variables

There is only one output format, listing the input data, observed and fitted count data (deaths or cases), coefficients, standard errors, relative risks (RR) and 95% confidence intervals, variance-covariance matrix, correlation matrix, goodness of fit statistics, and residuals for regression diagnostics. The program and all of its subroutines

are written in FORTRAN-77 and are run on a desktop PC with a AMD-K6-166MHz(MMX) chip. The algorithm can be compiled and run on Unix machines as well. The PC version of the executable requires 180,224 bytes of random access memory (RAM) to operate. Average execution time is several seconds per run, depending on the amount of input data.

2 PROGRAM DESCRIPTION

The computer program proceeds as follows. During run-time, the user interface with the program involves the following:

- Specifying the type of model to be fit (1-additive, 2-multiplicative, 3-power, and 4-non-linear)
- For additive models, specifying the power, ρ , which ranges from 1 for purely additive models to 0 for purely multiplicative models
- Specifying the number of independent variables (not cases or person-years) that are to be read from the input file
- Specifying the number of parameters to estimate
- For non-linear models, optional starting values for parameters
- For multiplicative models, specifying if the regression coefficients are to be exponentiated to obtain RR and its 95% CI
- Specifying whether or not the variance-covariance and correlation matrices are written to the output file whose name is specified by the user
- Specifying whether or not the input data are written to the output file whose name is specified by the user
- Specifying whether or not the observed and fitted count data (deaths or cases) are written to the output file
- Specifying whether or not the leverages, and standardized and deletion residuals are written to the output file

Once the data above have been specified during run time, the program does the following:

1. First, a “starting” regression is performed by regressing the ratio of counts (deaths or cases) to person-years of follow-up as the dependent variable on the independent variables. For non-linear models, the default starting regression can be skipped by specifying starting values for each parameter.
2. The difference between the observed and fitted counts (error residuals) from the starting regression are used as the dependent variable and regressed on the cross product of the inverse of the information matrix (variance covariance-matrix) and the score vector to obtain the solution vector.
3. The solution vector is then added to the original regression coefficients obtained during the starting regression.
4. Iterations are performed until the sum of the solution vector is below a fixed value called the *convergence criterion*, for which the default in PIRLS is 10^{-5} .
5. After convergence is reached, the standard errors are obtained from the diagonal of the variance-covariance matrix, and if specified, residuals for regression diagnostics are estimated.
6. Depending on specifications made by the user, various data are written to the output file, and then the program terminates run-time.

3 NOTATION AND THEORY

3.1 Binomial Basis of Poisson Regression

For large n and small p , e.g., a cancer rate of say 50/100,000 population, binomial probabilities are approximated by the Poisson distribution given by the form

$$P(x) = \frac{e^{-m}}{x!}, \quad (1)$$

Table 1: Deaths from coronary disease among British male doctors [21].

Age group, i	Non-smokers			Smokers		
	d_i	T_i	$\lambda_i(0)$	d_i	T_i	$\lambda_i(1)$
35-44	2	18,790	0.1064	32	52,407	0.6106
45-54	12	10,673	1.1243	104	43,248	2.4047
55-64	28	5,710	4.9037	206	28,612	7.1998
65-74	28	2,585	10.8317	186	12,663	14.6885
75-84	31	1,462	21.2038	102	5,317	19.1838

where the irrational number e is 2.71828.... In many respects, the Poisson distribution has been widely used in science and does not really have a direct relationship with the binomial distribution. As such, np is replaced by μ in the relationship

$$f(x) = \frac{\mu^x e^{-\mu}}{x!}, \quad (2)$$

where μ is the mean of the Poisson distribution. It has been said that as long as npq is less than 5, then the data are said to be Poisson distributed.

3.2 Poisson Assumption

The Poisson assumption states that, for stratified data, the number of deaths, d , in each cell approximates the values $x=0,1,2,\dots$ according to the formula

$$P(d = x) = \frac{(\lambda T)^x e^{-\lambda T}}{x!}, \quad (3)$$

where λ is the rate parameter which is equal to the number of deaths in each cell divided by the respective number of *person-years* (T) in that same cell. As an example, Table 1 lists a multiway contingency table for the British doctor data for coronary disease and smoking [21]. The rates for non-smokers, $\lambda_i(0)$, are simply the ratio of the number of deaths to T_i in age group i and exposure group 0, whereas the rates for smokers, $\lambda_i(1)$, are the ratio of deaths to T_i in age group i for the exposure group 1.

3.3 Ordinary Least Squares

Let us now look into the sum of squares for common linear regression models. We recall that most of our linear regression experience is with the model

$$\mathbf{Y} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad (4)$$

where \mathbf{Y} is a column vector of responses, \mathbf{X} is the data matrix, $\boldsymbol{\beta}$ is the column vector of coefficients and $\boldsymbol{\varepsilon}$ is the error. Another way of expressing the observed dependent variable, y_i , of (4) in *model* terms is

$$y_i = \beta_0 + \beta_1 x_{i1} + \beta_2 x_{i2} \cdots \beta_j x_{ij} + \varepsilon_i, \quad (5)$$

and in terms of the *predicted* values, \hat{y}_i , is

$$\hat{y}_i = \hat{\beta}_0 + \hat{\beta}_1 x_{i1} + \hat{\beta}_2 x_{i2} \cdots \hat{\beta}_j x_{ij} + e_i. \quad (6)$$

The residual sum of squares of (6) is

$$\begin{aligned} SS(\boldsymbol{\beta}) &= \sum_{i=1}^n \{y_i - \hat{y}_i\}^2 \\ &= \sum_{i=1}^n e_i^2 \\ &= \mathbf{e}'\mathbf{e}. \end{aligned} \quad (7)$$

Because of the relationship

$$\mathbf{e} = \mathbf{Y} - \mathbf{X}\boldsymbol{\beta}, \quad (8)$$

we can rewrite (7) in matrix notation as

$$SS(\boldsymbol{\beta}) = (\mathbf{Y} - \mathbf{X}\boldsymbol{\beta})'(\mathbf{Y} - \mathbf{X}\boldsymbol{\beta}). \quad (9)$$

3.4 Partitioning Error Sum of Squares for Weighted Least Squares

Suppose we have the general least squares equation

$$\mathbf{Y} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}. \quad (10)$$

and according to Draper and Smith [22] set

$$\mathbf{E}(\boldsymbol{\varepsilon}) = \mathbf{0} \quad \mathbf{V}(\boldsymbol{\varepsilon}) = \mathbf{V}\sigma^2 \quad \boldsymbol{\varepsilon} \sim \mathbf{N}(\mathbf{0}, \mathbf{V}\sigma^2), \quad (11)$$

and say that

$$\mathbf{P}'\mathbf{P} = \mathbf{P}\mathbf{P} = \mathbf{P}^2 = \mathbf{P}\mathbf{P} = \mathbf{V}. \quad (12)$$

Now if we let

$$\mathbf{f} = \mathbf{P}^{-1}\boldsymbol{\varepsilon} \quad \mathbf{E}(\mathbf{f}) = \mathbf{0}, \quad (13)$$

and then premultiply (10) by \mathbf{P}^{-1} , as in

$$\mathbf{P}^{-1}\mathbf{Y} = \mathbf{P}^{-1}\mathbf{X}\boldsymbol{\beta} + \mathbf{P}^{-1}\boldsymbol{\varepsilon}, \quad (14)$$

or

$$\mathbf{Z} = \mathbf{Q}\boldsymbol{\beta} + \mathbf{f}. \quad (15)$$

By applying basic least squares methods to (14) the residual sums of squares is

$$\mathbf{f}'\mathbf{f} = \boldsymbol{\varepsilon}'(\mathbf{P}^{-1})'\mathbf{P}^{-1}\boldsymbol{\varepsilon}, \quad (16)$$

since

$$(\mathbf{P}^{-1}\boldsymbol{\varepsilon})' = \boldsymbol{\varepsilon}'(\mathbf{P}^{-1})', \quad (17)$$

and since

$$(\mathbf{B}^{-1}\mathbf{A}^{-1}) = (\mathbf{A}\mathbf{B})^{-1}, \quad (18)$$

then

$$\mathbf{f}'\mathbf{f} = \boldsymbol{\varepsilon}'(\mathbf{P}\mathbf{P})^{-1}\boldsymbol{\varepsilon}, \quad (19)$$

which is equivalent to

$$\mathbf{f}'\mathbf{f} = \boldsymbol{\varepsilon}'(\mathbf{V})^{-1}\boldsymbol{\varepsilon}. \quad (20)$$

The sum of squares residual then becomes for (14)

$$SS(\boldsymbol{\beta}) = (\mathbf{Y} - \mathbf{X}\boldsymbol{\beta})'\mathbf{V}(\mathbf{Y} - \mathbf{X}\boldsymbol{\beta}). \quad (21)$$

3.5 Partitioning Error Sum of Squares for Iteratively Reweighted Least Squares

According to McCullagh and Nelder [23], a *generalized linear model* is composed of the following three parts:

- *Random* component, λ , which is normally distributed with mean $\frac{1}{T_i} \cdot \mu_i$ and variance $\frac{\mu_i}{T_i^2}$.
- *Systematic* component, x , which are covariates related to the linear predictor, $(\mathbf{x}_i'\boldsymbol{\beta})$, by the form

$$(\mathbf{x}_i'\boldsymbol{\beta}) = \sum_{j=1}^p x_{ij}\beta_j. \quad (22)$$

Table 2: The rate parameter and its various link functions in PIRLS.

Form of predictor for λ	Model	Risk coefficients
$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = (\mathbf{x}'_i\boldsymbol{\beta})$	Additive	Absolute Risk
$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = \exp(\mathbf{x}'_i\boldsymbol{\beta})$	Multiplicative	Relative Risk
$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = (\mathbf{x}'_i\boldsymbol{\beta})^p$	Power	N/A
$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = \beta_1 x_{i1} [1 - \{1 - e^{(-\beta_2 x_{i2})}\}^{\beta_3}]$	Non-linear	N/A

- The *link* between the random and systematic components for a linear model is given by

$$\lambda_i = \lambda_i(\mathbf{x}'_i\boldsymbol{\beta}), \quad (23)$$

where λ_i is the *observed* rate parameter and $\lambda_i = (\mathbf{x}'_i\boldsymbol{\beta})$ is the *predicted* rate parameter. It should be pointed out that the rate parameter, $\lambda_i = (\mathbf{x}'_i\boldsymbol{\beta})$, serves as its own link function. Thus, the rate parameter provides a linearization of our distributional models so that we can use multiple linear regression to obtain Best Linear Unbiased Estimators (BLUEs). In effect, the rate parameter allows us to reduce non-linear (non-Gaussian) models to linear models so that we can solve the system of linear equations. The rate parameter can take on various linear and non-linear functions. These are shown in Table 2. The power transformation was adapted from [24].

As shown in Table 2, it is obvious that the random component, λ_i , can take on a non-linear relationship with the systematic component, x . To *linearize* λ_i in a least squares approach by a succession of stages or *iterations*, we carry out a Taylor series expansion of λ_i about a point β_0 . Since we know that

$$g(\beta) = g(\beta_0) + g(\beta_0)'(\beta - \beta_0), \quad (24)$$

a Taylor series expansion of λ_i about $(\beta - \beta_0)$ gives us

$$\lambda_i \simeq \lambda_i(\mathbf{x}'_i\boldsymbol{\beta}_0) + \sum_{j=1}^p \left[\frac{\partial \lambda_i(\mathbf{x}'_i\boldsymbol{\beta})}{\partial \beta_j} \right]_{\boldsymbol{\beta} = \boldsymbol{\beta}_0} (\beta - \beta_0) + e_i. \quad (25)$$

Rearranging this equation, we have

$$\lambda_i - \lambda_i(\mathbf{x}'_i\boldsymbol{\beta}_0) = \mathbf{Z}^0 \boldsymbol{\delta}^0 + e_i, \quad (26)$$

where

$$\mathbf{Z}_{ij}^0 = \left[\frac{\partial \lambda_i(\mathbf{x}'_i\boldsymbol{\beta})}{\partial \beta_j} \right]_{\boldsymbol{\beta} = \boldsymbol{\beta}_0}, \quad (27)$$

and

$$\delta_i^0 = \beta_i - \beta_i^0. \quad (28)$$

The *Jacobian* matrix of partial derivatives of $\lambda_i(\mathbf{x}'_i\boldsymbol{\beta})$ with respect to the parameters, β at the zeroth iteration is then

$$\mathbf{Z}_0 = \begin{bmatrix} z_{11} & z_{12} & \cdots & z_{1p} \\ z_{21} & z_{22} & \cdots & z_{2p} \\ \vdots & \vdots & \ddots & \vdots \\ z_{n1} & z_{n2} & \cdots & z_{np} \end{bmatrix} = \mathbf{Z}_{ij}^0, n \times p. \quad (29)$$

From (26), the residuals, e_i , now become

$$e_i = \lambda_i - \hat{\lambda}_i(\mathbf{x}'_i\boldsymbol{\beta}) - \mathbf{Z}\boldsymbol{\delta}. \quad (30)$$

3.5.1 Development of Dependent Variable, \mathbf{Y}

Using (30), let $\lambda_i - \lambda_i^0(\mathbf{x}_i' \boldsymbol{\beta})$ be equal to Y_i . We now rewrite the residuals

$$e_i = Y_i - \mathbf{Z}\boldsymbol{\delta}, \quad (31)$$

so that the column vector, \mathbf{Y} of dependent \mathbf{Y} becomes

$$\mathbf{Y} = \begin{bmatrix} \lambda_1 - \lambda_1^0(\mathbf{x}_1' \boldsymbol{\beta}) \\ \lambda_2 - \lambda_2^0(\mathbf{x}_2' \boldsymbol{\beta}) \\ \vdots \\ \lambda_n - \lambda_n^0(\mathbf{x}_n' \boldsymbol{\beta}) \end{bmatrix}. \quad (32)$$

3.5.2 Development of Weighting Matrix, \mathbf{W}

We know that $\text{Var}(cX) = c^2 \text{Var}(X)$ so if we substitute in $\hat{\lambda}$ we have

$$\begin{aligned} \text{Var}\left(\frac{1}{T_i} \cdot \hat{\mu}_i\right) &= \left(\frac{1}{T_i}\right)^2 \text{Var}(\hat{\mu}_i) \\ &= \frac{\hat{\mu}_i^2}{T_i^2}. \end{aligned} \quad (33)$$

Therefore, each diagonal element of the *weight* matrix becomes

$$w_{ii} = \frac{1}{\text{Var}(\hat{\lambda}_i)} = \frac{T_i^2}{\hat{\mu}_i^2} = \frac{T_i \cdot T_i}{\hat{\lambda}_i \cdot T_i} = \frac{T_i}{\hat{\lambda}_i}. \quad (34)$$

Because λ_i can be very small and T_i very large and it is better to divide by a larger number than a smaller one with a computer, we will choose the weight $\frac{T_i^2}{\hat{\mu}_i^2}$. The entire weighting matrix \mathbf{W} is then

$$\mathbf{W} = \begin{bmatrix} \frac{T_1^2}{\hat{\mu}_1^2} & 0 & 0 & 0 \\ 0 & \frac{T_2^2}{\hat{\mu}_2^2} & 0 & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & \frac{T_n^2}{\hat{\mu}_n^2} \end{bmatrix}. \quad (35)$$

3.5.3 Defining Maximum Likelihood Solution Vector, $\boldsymbol{\delta}$

Let $\mathbf{V}^{-1} = \mathbf{W}$, and now substitute our new definition for residuals from (31) into (21) obtain the new sum of squares

$$SS(\boldsymbol{\delta}) = (\mathbf{Y} - \mathbf{Z}\boldsymbol{\delta})' \mathbf{W} (\mathbf{Y} - \mathbf{Z}\boldsymbol{\delta}) \quad (36)$$

$$= \mathbf{W}\mathbf{Y}'\mathbf{Y} - (\mathbf{Z}\boldsymbol{\delta})' \mathbf{W} \mathbf{Y} - \mathbf{Y}' \mathbf{Z} \mathbf{W} \boldsymbol{\delta} + \boldsymbol{\delta}' \mathbf{Z}' \mathbf{W} \mathbf{Z} \boldsymbol{\delta}. \quad (37)$$

The transpose $(\mathbf{Z}\boldsymbol{\delta})'$ of the term $(\mathbf{Z}\boldsymbol{\delta})' \mathbf{W} \mathbf{Y}$ in (37) is equal to $\boldsymbol{\delta}' \mathbf{Z}'$ because $(\mathbf{A}\mathbf{B})' = \mathbf{B}' \mathbf{A}'$. In addition, since $(\mathbf{Z}\boldsymbol{\delta})' \mathbf{W} \mathbf{Y}$ is a 1×1 matrix, or scalar, then its transpose $\mathbf{Y}' \mathbf{Z} \mathbf{W} \boldsymbol{\delta}$ has the same value. Thus, (37) is transformed into

$$SS(\boldsymbol{\delta}) = \mathbf{W}\mathbf{Y}'\mathbf{Y} - 2\boldsymbol{\delta}' \mathbf{Z}' \mathbf{W} \mathbf{Y} + \boldsymbol{\delta}' \mathbf{Z}' \mathbf{W} \mathbf{Z} \boldsymbol{\delta}. \quad (38)$$

The $\boldsymbol{\delta}'$ on the left sides of parameters in the right side of (38) drop out because if we set

$$\mathbf{A} = (\mathbf{Z}' \mathbf{W} \mathbf{Y}), \quad (39)$$

and from Searle [25] know that

$$\frac{\partial \mathbf{x}' \mathbf{A}}{\partial \mathbf{x}} = \mathbf{A}, \quad (40)$$

and if we set

$$\mathbf{A} = (\mathbf{Z}'\mathbf{W}\mathbf{Z}\boldsymbol{\delta}), \quad (41)$$

and from [25] know that

$$\frac{\partial(\mathbf{x}'\mathbf{A}\mathbf{x})}{\partial\mathbf{x}} = 2\mathbf{A}\mathbf{x}, \quad (42)$$

then the normal equation then becomes

$$SS(\boldsymbol{\delta}) = \mathbf{Z}'\mathbf{W}\mathbf{Z}\boldsymbol{\delta} + \mathbf{Z}'\mathbf{W}\mathbf{Y}. \quad (43)$$

In order to minimize (43) with respect to the $\boldsymbol{\delta}$ parameters, we must take the first partial derivative. This gives us

$$\frac{\partial SS(\boldsymbol{\delta})}{\partial\boldsymbol{\delta}} = -2\mathbf{Z}'\mathbf{W}\mathbf{Y} + 2(\mathbf{Z}'\mathbf{W}\mathbf{Z}\boldsymbol{\delta}), \quad (44)$$

that provides us with the consistent equations

$$(\mathbf{Z}'\mathbf{W}\mathbf{Z}\boldsymbol{\delta}) = \mathbf{Z}'\mathbf{W}\mathbf{Y}, \quad (45)$$

that has solution

$$\boldsymbol{\delta} = (\mathbf{Z}'\mathbf{W}\mathbf{Z})^{-1}\mathbf{Z}'\mathbf{W}\mathbf{Y}. \quad (46)$$

At the i th iteration, the values of the parameters are

$$\boldsymbol{\beta}_i = \boldsymbol{\beta}_{i-1} + \boldsymbol{\delta}. \quad (47)$$

Convergence is reached at the point when

$$\{\boldsymbol{\beta}_i - \boldsymbol{\beta}_{(i-1)}\}/\boldsymbol{\beta}_{(i-1)} < \epsilon, \quad (48)$$

where ϵ is 10^{-5} by default in PIRLS.

3.6 Variance-covariance Matrix

The variance-covariance matrix, with parameter variance on the diagonal and covariances on the off-diagonals, at convergence is identical to the information matrix or inverse of the weighted dispersion matrix

$$\mathbf{I}_{\boldsymbol{\beta}}^{-1} = (\mathbf{Z}'\mathbf{W}\mathbf{Z})^{-1} = \begin{bmatrix} \sigma_1^2 & \sigma_{12} & \cdots & \sigma_{1j} \\ \sigma_{12} & \sigma_2^2 & \cdots & \sigma_{2j} \\ \vdots & \vdots & \ddots & \vdots \\ \sigma_{1j} & \sigma_{2j} & \cdots & \sigma_j^2 \end{bmatrix}. \quad (49)$$

3.7 Standard Errors of Regression Coefficients

The standard errors of the regression coefficients are equivalent to $\sqrt{\sigma_j^2}$. A Wald test for each parameter is simply $\beta_j/\text{s.e.}(\beta_j)$ and indicates significance if the ratio is < -1.96 or > 1.96 at the $\alpha = 0.05$ level of significance.

3.8 Residuals and Regression Diagnostics

The error residual, e_i has mean 0 and measures the lack of concordance of the observed μ_i against the fitted value $\hat{\mu}_i$. The greater the value of e_i , the worse the fit of the model. Pearson residuals, r_P are estimate as

$$r_P = \frac{\mu_i - \hat{\mu}_i}{\mu}, \quad (50)$$

and provide a measure of the residual variation. The deviance residuals

$$r_D = \mu_i \left[\log \frac{\mu_i}{\hat{\mu}_i} \right] + (\hat{\mu}_i - \mu_i), \quad (51)$$

allow the investigator to determine the goodness of fit of the model, i.e., how well the observed data agree with the fitted values. Freeman and Tukey [26] introduced a variance stabilized residual for Poisson models given by

$$r_{FT} = \sqrt{\mu} + \sqrt{\mu + 1} - \sqrt{\hat{\mu} + 1}. \quad (52)$$

The leverage [23] for each observation, h_i , provides information about the geometric distance between a given point $(z_{i1}, z_{i2}, \dots, z_{ij})$ and the centroid of all points $(\bar{z}_{i1}, \bar{z}_{i2}, \dots, \bar{z}_{ij})$ in the predictor space. Individual leverages, h_i , are obtained from the diagonal of the “Hat” matrix given by

$$\mathbf{H} = \sqrt{\mathbf{W}}\mathbf{Z}(\mathbf{Z}'\mathbf{W}\mathbf{Z})^{-1}\mathbf{Z}'\sqrt{\mathbf{W}}. \quad (53)$$

Because $\sum h_i = \text{trace}(\mathbf{H}) = p$, observations with high leverage can approach unity. A criterion introduced by Hoaglin and Welsch [27] for determining high leverage for an observation states that a point has high leverage if $h_i > 2p/n$. Points with high leverage are located at a substantial distance from the fitted regression line, and act to “pull” the fitted line toward their location. Increases in the goodness of fit of a model can be obtained by refitting a model after the observation(s) with high leverage have been removed.

Cook [28] introduced the residual, $\Delta(\beta)_{-ij}$, which measures the difference between β when the observation is included in the model and β when the observation is not in the model. The more positive the values of $\Delta(\beta)_{-ij}$, the more influence an observation has when compared with other observations. Individual $\Delta(\beta)_{-ij}$ residuals are obtained by the $n \times p$ matrix $\Delta(\beta)_{-ij}$, by the relationship

$$\Delta(\beta)_{-ij} = \frac{-(\mathbf{Z}'\mathbf{W}\mathbf{Z})^{-1}z_{ij}e_{ij}w_{ij}}{1 - h_{ij}}. \quad (54)$$

Finally, the *standardized* residual is used to measure the model error introduced by particular observation when its error residual has constant variance

$$\tilde{r}_i = \frac{e_i}{\sqrt{(1 - h_i)}}. \quad (55)$$

Values for \tilde{r}_i that exceed 1.96 tend to be outliers at the $\alpha = 0.05$ level of significance.

3.9 Goodness-of-Fit Statistics

The Pearson χ^2 goodness of fit is

$$\chi^2 = \sum_{i=1}^n r_P^2, \quad (56)$$

and the deviance goodness of fit is

$$D = \sum_{i=1}^n r_D, \quad (57)$$

which is also χ^2 distributed. If a model fits, its χ^2 and D will be lower than the degrees of freedom $(n - p)$.

3.10 Structure

Program flow is outline in Figure 1. The starting regression to obtain the initial values of the parameter vector, β_i is performed by calling:

CALL INIT(NREG, NCOL, NROW, ISTART, START, B, G, RHO, IPRED, IDATA, INFILE, OUTFILE, IFLT)

For non-linear models, the user can specify starting values for the parameters and skip the multiplicative fit to obtain starting parameter estimates. The solution vector, δ , used to modify the original regression coefficients during each iteration are determined through calls of:

CALL PRED(NCOL, NROW, G, IDF, DEV, TPR, NREG, RHO, IPRED, ZTWZ1, ZTWY, ITER, WIN, YDIFF, OEXPE, OOPS, OFT, OPR, XIN, XT, IFLT)

Lastly, the leverages, deletion residuals, and standardized residuals are obtained by calling:

CALL FIT(WIN, YDIFF, XIN, XT, ZTWZ1, OEXPE, LEV, DRES, SRES, NCOL, NROW, IFLT)

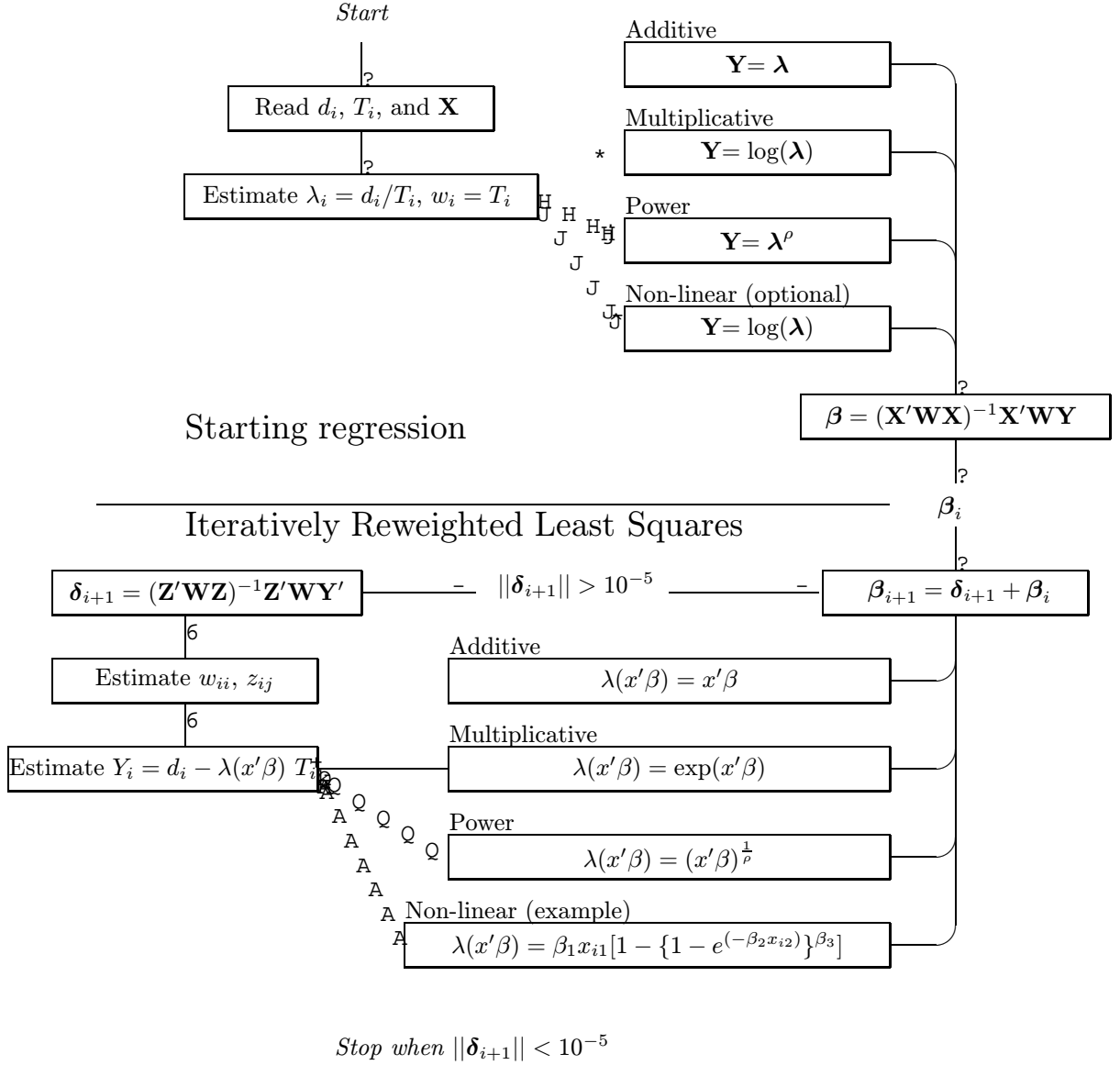


Figure 1: Program flow in PIRLS.

The parameter, arrays, and variables used in PIRLS are defined below.

MP=4000	Parameter	Constant	Fixed array size for maximum number of records
NP=28	Parameter	Constant	Fixed array size for maximum number of variables
IP=1	Parameter	Constant	Fixed array size for when 1 is needed
NIP=16	Parameter	Constant	Fixed array when more than 1 is needed
INFILE	Character	Input	Input filename and path
OUTFILE	Character	Input	Output filename and path
NROW	Integer	Input	Number of input records read
NCOL	Integer	Input	Number of parameters specified by user
ISTART	Integer	Input	For non-linear models; 0=user specified starting values, 1=starting values estimated by multiplicative fit in subroutine INIT
START	Real	Input	For non-linear models; starting values specified by user when ISTART=0
IPRED	Integer	Input	Number of parameters to be read in from input file
NREG	Integer	Input	Type of regression model to be fit (1-additive, 2-multiplicative, 3-power, 4-non-linear)
IDATA	Integer	Input	Write input data to output file, 1=yes, 0=no
ICOVAR	Integer	Input	Write variance-covariance and correlation matrices to output file, 1=yes, 0=no
ICELLS	Integer	Input	Write observed and fitted counts, d_i , to output file, 1=yes, 0=no
IEX	Integer	Input	Exponentiate regression coefficients for multiplicative models only, i.e., NREG=2
IREG	Integer	Input	Estimate standardized and deletion residuals and leverages, 1=yes, 0=no
RHO	Real	Input	Power transform for power models, ρ
OBS(MP)	Real	Input	Number of deaths or cases in subgroup (record), μ_i

T(MP)	Real	Input	Number of person-years of follow-up in subgroup (record), T
X(MP,NP)	Real	Work	Design matrix of input independent variables, \mathbf{X}
LAMBDA(MP)	Real	Work	Observed rate, λ_i
LAMHAT(MP)	Real	Work	Predicted rate, $\hat{\lambda}_i$
EXPE(MP)	Real	Work	Expected deaths or cases in subgroup, $\hat{\mu}_i = \hat{\lambda}_i \times T$
Y(MP)	Real	Work	Error Residual, $Y_i = \mu_i - \hat{d}_i$
W(MP)	Real	Work	Variance weights = $\mathbf{W} = T^2 / \text{Var}(\hat{\lambda}_i)$
Z(MP,NP)	Real	Work	Jacobian matrix of first partial derivatives, \mathbf{Z}
ZTWZ(MP,NP)	Real	Work	Information matrix, $(\mathbf{Z}'\mathbf{W}\mathbf{Z})$
ZTWZ1(NP,NP)	Real	Work	Inverse of information matrix, $(\mathbf{Z}'\mathbf{W}\mathbf{Z})^{-1}$
ZTWY(NP,1)	Real	Work	Score vector, $\mathbf{Z}'\mathbf{W}\mathbf{Y}$
FT(MP)	Real	Work	Freeman-Tukey residuals, r_{FT}
PR(MP)	Real	Work	Pearson χ^2 residuals, r_P
DEV	Real	Work	Deviance residuals, r_D
VARCOV(NP,NP)	Real	Work	Variance-covariance matrix used for output, $(\mathbf{Z}'\mathbf{W}\mathbf{Z})$
G(NP,1)	Real	Work	Solution vectors, δ
B(NP,1)	Real	Work	Regression coefficients updated at each iteration, β_i
SE(NP)	Real	Work	Standard error of β_i
CORR(NP,NP)	Real	Work	Correlation matrix of coefficients
LL(NP)	Real	Work	Lower bound of 95% confidence limit for relative risk (multiplicative model only)
UL(NP)	Real	Work	Upper bound of 95% confidence limit for relative risk (multiplicative model only)
ML(NP)	Real	Work	Relative risk (multiplicative model only)
LEV(MP,MP)	Real	Work	Leverages from the Hat matrix, h_i
DRES(MP,NP)	Real	Work	Deletion residuals, $\Delta(\beta)_{i,j}$
SRES(MP)	Real	Work	Standardized residuals, \tilde{r}_i
IDF	Integer	Work	Degrees of freedom used for goodness of fit
DEV	Real	Work	Deviance goodness of fit, $D_{d.f.}^2$
TPR	Real	Work	Pearson χ^2 goodness of fit $\chi_{d.f.}^2$
IFLT	Integer	Output	Premature exit codes 0=Convergence reached 1=NCOL < 1 2=NROW < 1 3=Degrees of freedom, (NROW-NCOL), < 1 4= $ \sum \beta_j = 0$ 5= $ \sum G_j = 0$ 6=Iterations > 100

3.11 Auxiliary Algorithms

MATMUL [29] performs matrix multiplication for the matrix manipulation. *TRNPOS* transposes arrays as needed for matrix multiplication. *MATINV* calls *SVDCMP* based on [30] to invert the information matrix using singular value decomposition. *ALG* and *GAMAIN* are used for estimating tabled values of χ^2 .

3.12 Restriction

All input parameters are checked for nullity and a fault message is returned if there is an illegal entry.

3.13 Time

A thorough investigation of absolute timing has not been performed; however, it should be noted that execution time is a function of the number of model parameters and effects to be considered.

3.14 Precision

The algorithm may be converted to double precision by making the following changes:

1. Change *REAL* to *DOUBLE PRECISION* in the algorithm.
2. Change the constants to double precision.
3. Change *EXP* to *DEXP* and *ALOG* to *DLOG* in all applicable routines.
4. Make appropriate changes in auxiliary routines *MATMUL*, *TRNPOS*, *MATINV*, *SVDCMP*, and in functions *ALG* and *GAMAIN*.

3.15 Application - Fitting a Multiplicative Model

3.15.1 Data Arrangement and Partial Derivatives

We shall enumerate the steps required for fitting a multiplicative model. Before we start the regression run, however, we must set up the data file and also determine the first partial derivatives of the rate ratio λ with respect to each parameter in the model.

1. **Arranging the data in an ASCII text file.** The first step is to arrange our data in an ASCII text file. The typical ASCII text file arrangement of data for a Poisson regression run using data in Table 1 is shown in the following:

Cases	T/1,000	Age	Smoke
2.	18.790	1	0 0 0 0 0
12.	10.673	0	1 0 0 0 0
28.	5.710	0	0 1 0 0 0
28.	2.585	0	0 0 1 0 0
31.	1.462	0	0 0 0 1 0
32.	52.407	1	0 0 0 0 1
104.	43.248	0	1 0 0 0 1
206.	28.612	0	0 1 0 0 1
186.	12.663	0	0 0 1 0 1
102.	5.317	0	0 0 0 1 1

where the person-years in Table 1 are divided by 1,000 and the covariates denoting the age group are dummy coded (0,1) with a 1 to indicate that the rates are from a given cell, and whether or not the subgroup smoked.

2. **Determining Partial Derivatives of $\lambda_i(\mathbf{x}'_i|\boldsymbol{\beta})$ With Respect to Each β_j .** Next we must write out the partial derivatives of the rate parameter with respect to each parameter (these will be used later in the algorithm). The partial derivatives are used in the Jacobian \mathbf{Z} matrix to form the scores during each iteration. Thus, each partial derivative forms an element, z_{ij} , in the Jacobian \mathbf{Z} matrix. Given that the partial derivative of an exponentiated function u in the form e^u is $e^u du$, the partial derivative of $\lambda_i(\mathbf{x}'_i|\boldsymbol{\beta})$ with respect to each parameter, β_j is as follows:

$$z_{ij} = \frac{\partial \lambda_i(\mathbf{x}'_i|\boldsymbol{\beta})}{\partial \beta_j} x_{ij}, \quad (58)$$

3.15.2 Steps in the Algorithm

The following section discusses the steps taken by the algorithm for performing iteratively reweighted least squares.

1. **Definition of the model.** The model we will fit is a log-linear model

$$\lambda_i(\mathbf{x}'_i|\boldsymbol{\beta}) = e^{(\beta_1 x_{i1} + \beta_2 x_{i2} + \beta_3 x_{i3} + \beta_4 x_{i4} + \beta_5 x_{i5} + \beta_6 x_{i6})}, \quad (59)$$

which is equivalent to

$$\log(\lambda_i(\mathbf{x}'_i|\boldsymbol{\beta})) = (\beta_1 x_{i1} + \beta_2 x_{i2} + \beta_3 x_{i3} + \beta_4 x_{i4} + \beta_5 x_{i5} + \beta_6 x_{i6}). \quad (60)$$

2. **Estimate column vector of dependent variables for the log-linear model.** The algorithm starts by reading in the data from the input file and then estimates the rate parameter, λ_i , for each record as the ratio of number of *observed* deaths to the number of *observed* person-years. The log of λ_i is then calculated to obtain the dependent variable column vector

$$\mathbf{Y} = \begin{bmatrix} \log\left(\frac{d_1}{T/1,000}\right) \\ \log\left(\frac{d_2}{T/1,000}\right) \\ \log\left(\frac{d_3}{T/1,000}\right) \\ \log\left(\frac{d_4}{T/1,000}\right) \\ \log\left(\frac{d_5}{T/1,000}\right) \\ \log\left(\frac{d_6}{T/1,000}\right) \\ \log\left(\frac{d_7}{T/1,000}\right) \\ \log\left(\frac{d_8}{T/1,000}\right) \\ \log\left(\frac{d_9}{T/1,000}\right) \\ \log\left(\frac{d_{10}}{T/1,000}\right) \end{bmatrix}. \quad (61)$$

and when substituting in the data above we get

$$\mathbf{Y} = \begin{bmatrix} \log\left(\frac{2}{18,790}\right) \\ \log\left(\frac{12}{10,673}\right) \\ \log\left(\frac{28}{5,710}\right) \\ \log\left(\frac{28}{2,585}\right) \\ \log\left(\frac{31}{1,462}\right) \\ \log\left(\frac{32}{52,407}\right) \\ \log\left(\frac{104}{43,248}\right) \\ \log\left(\frac{206}{28,612}\right) \\ \log\left(\frac{186}{12,663}\right) \\ \log\left(\frac{102}{5,317}\right) \end{bmatrix}. \quad (62)$$

3. **Construction of the \mathbf{X} matrix of independent variables.** In addition to the estimation of the dependent variable \mathbf{Y} column vector, the algorithm constructs the \mathbf{X} matrix of independent variables from the independent variables read in from the input file as

$$\mathbf{X} = \begin{bmatrix} 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 1 & 0 & 0 & 0 & 0 & 1 \\ 0 & 1 & 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 & 0 & 1 \\ 0 & 0 & 0 & 1 & 0 & 1 \\ 0 & 0 & 0 & 0 & 1 & 1 \end{bmatrix}. \quad (63)$$

4. **“Starting Regression” - log-linear regression to estimate initial values of parameters.** Once the \mathbf{Y} and \mathbf{X} matrices are constructed, the algorithm fits the log-linear model

$$\log(\lambda_i(\mathbf{x}'_i\boldsymbol{\beta})) = (\beta_1x_1 + \beta_2x_2 + \beta_3x_3 + \beta_4x_4 + \beta_5x_5 + \beta_6x_6), \quad (64)$$

which in matrix terms is given as

$$\boldsymbol{\beta} = (\mathbf{X}'\mathbf{W}\mathbf{X})^{-1}\mathbf{X}'\mathbf{W}\mathbf{Y}, \quad (65)$$

where diagonal elements of \mathbf{W} are set equal to T . Figure 1 shows the program flow in PIRLS for the various models. Table 3 shows the link functions for λ for each starting regression and the linear predictor used for estimating $\lambda_i(\mathbf{x}'_i\boldsymbol{\beta})$ during subsequent iterations.

Table 3: Link functions used during “starting regression” and linear predictor used to estimate $\hat{\lambda}$ during subsequent iterations.

Model	Link function for λ	Form of predictor for $\lambda_i(\mathbf{x}'_i\boldsymbol{\beta})$
Additive	λ_i	$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = (\mathbf{x}'_i\boldsymbol{\beta})$
Multiplicative	$\log(\lambda_i)$	$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = \exp(\mathbf{x}'_i\boldsymbol{\beta})$
Power	λ_i^p	$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = (\mathbf{x}'_i\boldsymbol{\beta})^{\frac{1}{p}}$
Non-linear (optional)*	$\log(\lambda_i)$	$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = \exp(\mathbf{x}'_i\boldsymbol{\beta})$

* For non-linear models, the user can specify starting values for parameters, rather than using the default multiplicative fit to obtain starting values.

5. **Estimation of predicted rate parameter.** After the first regression or **iteration**, the regression coefficients, β_j , are used to estimate predicted rates from the linear predictor

$$\lambda_i(\mathbf{x}'_i\boldsymbol{\beta}) = e^{\sum_{j=1}^p x_{ij}\beta_j} = e^{(\mathbf{x}'_i\boldsymbol{\beta})}. \quad (66)$$

6. **Estimation of expected number of deaths.** The predicted rates are then multiplied by the observed person-years in each record to give the expected number of deaths as

$$\hat{\mu}_i = d_i(\mathbf{x}'_i\boldsymbol{\beta}) = \lambda_i(\mathbf{x}'_i\boldsymbol{\beta})T_i. \quad (67)$$

7. **Estimation of dependent variable.** The expected number of deaths, $\hat{\mu}$, are then subtracted from the observed deaths, μ , to give the dependent variable

$$Y_i = \mu_i - \hat{\mu}_i = d_i - \lambda_i(\mathbf{x}'_i\boldsymbol{\beta})T_i. \quad (68)$$

8. **Construction of column vector \mathbf{Y} .** The column vector \mathbf{Y} is then equal to $[Y_i]$.

Table 4: Survival data for splenic stem cell colony formation after irradiation and transplantation (see ref. [31]).

Colonies, μ	Trials, T_i	Cell concentration, x_{i1}	Dose(Sv), x_{i2}
60.0	6.0	1.25	0.00
66.0	7.0	1.75	0.96
46.0	4.0	3.00	1.92
82.0	9.0	7.20	2.88
105.0	11.0	24.00	4.32
123.0	15.0	75.00	5.76
12.0	4.0	120.00	6.72

9. **Estimation of the partial derivatives for each row vector of the data matrix \mathbf{X} .** For each record in the data matrix, \mathbf{X} , estimate the value of the partial derivatives using the current parameter values, β_j . For our log-linear model in (65), we get

$$z_{ij} = \frac{\partial \lambda_i(\mathbf{x}'_i \boldsymbol{\beta})}{\partial \beta_j} = e^{(\mathbf{x}'_i \boldsymbol{\beta})} x_{ij}. \quad (69)$$

10. **Construction of the Jacobian matrix of partial derivatives.** Now the algorithm constructs the $n \times p$ Jacobian matrix \mathbf{Z} composed of the first partial derivatives estimated above, which appears as $\mathbf{Z} = [z_{ij}]$.
11. **Construction of the inverse variance weighting matrix.** Since the Poisson variance of $\hat{\mu}_i$ changes at each iteration, we must weight the dispersion matrix with weights defined by $T_i^2 / \hat{\mu}_i$ given as $\mathbf{W} = \text{diag}[w_{ii}]$, where w_{ii} is $\frac{T_i^2}{\hat{\mu}_i}$.
12. **Estimation the solution vector at each iteration.** The maximum likelihood solutions at the next and subsequent iterations are obtained through the matrix manipulation

$$\boldsymbol{\delta} = (\mathbf{Z}' \mathbf{W} \mathbf{Z})^{-1} \mathbf{Z}' \mathbf{W} \mathbf{Y}. \quad (70)$$

13. **Updating regression coefficients.** At each iteration, the new solution vectors are added to the previous parameter estimates given as

$$\boldsymbol{\beta}_i = \boldsymbol{\beta}_{i-1} + \boldsymbol{\delta}. \quad (71)$$

14. **Repeat steps 5-13 above until convergence is reached.** The above steps (5 through 13) are repeated until the euclidean norm of the score vector is below some non-negative value

$$\{\beta_i - \beta_{(i-1)}\} / \beta_{(i-1)} < \epsilon, \quad (72)$$

where ϵ is 10^{-5} by default in PIRLS. At this point, convergence is reached at a global maximum.

3.16 Application - Fitting a Non-linear Model

For this example, we will fit a non-linear model to estimate the number of stem cell colonies in spleens of laboratory mice into which irradiated cells were transplanted [31]. Table 4 shows the colonies formed, μ , number of trials, T_i , cell concentration, x_{i1} , and radiation dose (Sv), x_{i2} , arranged in tabular notation.

The non-linear model we will fit to obtain maximum likelihood estimates of colonies formed is of the form

$$\lambda_i = \beta_1 x_{i1} [1 - \{1 - e^{(-\beta_2 x_{i2})}\}^{\beta_3}], \quad (73)$$

where x_{i1} is the concentration of transplanted cells, x_{i2} is the radiation dose (Sv), and β_1 , β_2 , and β_3 are the parameters of interest.

In order to build the Jacobian matrix, \mathbf{Z} , we need to take the first partial derivatives of λ_i with respect to β_1 , β_2 , and β_3 , i.e., $\partial\lambda_i/\partial\beta_1$, $\partial\lambda_i/\partial\beta_2$, and $\partial\lambda_i/\partial\beta_3$. Since there are three β coefficients, the first column entries for row i of \mathbf{Z} , i.e., z_{i1} , will be estimated by use of equation

$$\frac{\partial\lambda_i}{\partial\beta_1} = x_{i1}[1 - \{1 - e^{(-\beta_2 x_{i2})}\}^{\beta_3}]. \quad (74)$$

The first partial derivative in the 2nd column entry of row i of \mathbf{Z} , i.e., z_{i2} is estimated as

$$\frac{\partial\lambda_i}{\partial\beta_2} = -\beta_1 x_{i1} \{1 - e^{(-\beta_2 x_{i2})}\}^{\beta_3} \beta_3 x_{i2} \frac{e^{-\beta_2 x_{i2}}}{(1 - e^{-\beta_2 x_{i2}})}. \quad (75)$$

The third column of \mathbf{Z} has elements z_{i3} , which are numerically estimated as

$$\frac{\partial\lambda_i}{\partial\beta_3} = -\beta_1 x_{i1} \{1 - e^{(-\beta_2 x_{i2})}\}^{\beta_3} \ln(1 - e^{-\beta_2 x_{i2}}). \quad (76)$$

When fitting this model, we chose to specify starting values of $\beta_1 = 7.6364$, $\beta_2 = 0.9341$, and $\beta_3 = 7.6364$ from [3], rather than use the default multiplicative fit to obtain starting values. Input and output for this model using the data in Table 4 are shown in the Appendix.

3.17 Input

During run-time, the user must respond to the following queries concerning each run:

- The type of model to be fit (1-additive, 2-multiplicative, 3-power, and 4-non-linear)
- For additive models, the power, ρ , which ranges from 1 for purely additive models to 0 for purely multiplicative models
- The number of independent (input) variables to be read in from the input file
- The number of parameters to estimate
- For non-linear models, optional starting values for parameters
- For multiplicative models, the regression coefficients are to be exponentiated to obtain RR and its 95% CI
- Whether or not the variance-covariance and correlation matrices are written to the output file
- Whether or not the input data are written to the output file
- Whether or not the observed and fitted count data (deaths or cases) are written to the output file
- Whether or not the leverages, and standardized and deletion residuals are written to the output file

4 USER INPUT DATA FILE

PIRLS uses one input data file (ASCII) whose name is specified by the user at run-time. The input file contains three types of information

1. Run title (text) in record 1
2. FORTRAN format statement (in parentheses) in record 2 needed for reading the input data in later records
3. Input data, i.e., cases, person-years, and independent variables in all subsequent records

The examples in the Appendix give listings of input files.

5 SAMPLE INPUT/OUTPUT CASES

The output for all runs is in tabular form in the ASCII text file whose name is specified by the user at run-time. Four examples using an input file are given in the Appendix.

Table 5: Names and descriptions of each PIRLS file.

Filename	Description	Format
INFILE*	Input filename	ACSII text
OUTFILE*	Output filename	ACSII text
PIRLS.FOR	FORTTRAN source code	ASCII text
PIRLS.EXE	Executable file	FORTTRAN

* Input and output filenames are specified by user at run-time.

6 FILENAMES

Table 5 list the names of the files that were used to program, link and execute PIRLS. As one notices, the source code is an ASCII text file and the object and executable files have been compiled with Microsoft FORTRAN Powerstation 4.0. There is one required input data file, whose name is specified at run-time.

7 AVAILABILITY

The program and all of its subroutines are available from the Journal of Statistical Software free of charge at <http://www.stat.ucla.edu/journals/jss/>

8 ACKNOWLEDGEMENTS

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9 APPENDIX: SAMPLE INPUT/OUTPUT CASES

Example 1: Additive (linear) model using data in Table 1 adapted from [21]. The input data file for this run is distribution file EXAMP1-3.DAT and is listed below:

```
PIRLS
(F4.0,F8.3,6F3.0)
  2.  18.790 1. 0. 0. 0. 0. 0.
 12.  10.673 0. 1. 0. 0. 0. 0.
 28.   5.710 0. 0. 1. 0. 0. 0.
 28.   2.585 0. 0. 0. 1. 0. 0.
 31.   1.462 0. 0. 0. 0. 1. 0.
 32.  52.407 1. 0. 0. 0. 0. 1.
104.  43.248 0. 1. 0. 0. 0. 1.
206.  28.612 0. 0. 1. 0. 0. 1.
186.  12.663 0. 0. 0. 1. 0. 1.
102.   5.317 0. 0. 0. 0. 1. 1.
```

The screen queries to be entered by the user are as follows:

```
ENTER THE NAME OF INPUT FILE:
EXAMP1-3.DAT
```

```
ENTER THE NAME OF INPUT FILE:
EXAMPLE1.OUT
```

```
ENTER 1-ADD 2-MULT 3-POWER 4-NON-LINEAR
1
```

```
ENTER RHO:  1 - ADDITIVE      - OTHER
1
```

```
ENTER NUMBER OF PREDICTOR (INPUT) VARIABLES
6
```

```
NUMBER OF PARAMETERS TO BE FITTED?
6
```

```
PRINT INFORMATION AND CORR MATRICES?  0-NO    1-YES
1
```

```
PRINT INPUT DATA?  0-NO    1-YES
1
```

```
PRINT FITTED CELLS?  0-NO    1-YES
1
```

```
PERFORM REGRESSION DIAGNOSTICS?  0-NO    1-YES
1
```

The output file contains:

```
TITLE: PIRLS
```

MODEL SELECTED: ADDITIVE

INPUT DATA FORMAT: (F4.0,F8.3,6F3.0)

```

2.  18.790 1. 0. 0. 0. 0. 0.
12.  10.673 0. 1. 0. 0. 0. 0.
28.   5.710 0. 0. 1. 0. 0. 0.
28.   2.585 0. 0. 0. 1. 0. 0.
31.   1.462 0. 0. 0. 0. 1. 0.
32.  52.407 1. 0. 0. 0. 0. 1.
104. 43.248 0. 1. 0. 0. 0. 1.
206. 28.612 0. 0. 1. 0. 0. 1.
186. 12.663 0. 0. 0. 1. 0. 1.
102.   5.317 0. 0. 0. 0. 1. 1.

```

ITERATION:	1	DEVIANCE:	86.7977
ITERATION:	2	DEVIANCE:	89.7618
ITERATION:	3	DEVIANCE:	93.6565
ITERATION:	4	DEVIANCE:	98.2919
ITERATION:	5	DEVIANCE:	22.3763
ITERATION:	6	DEVIANCE:	7.7339
ITERATION:	7	DEVIANCE:	7.4384
ITERATION:	8	DEVIANCE:	7.4331
ITERATION:	9	DEVIANCE:	7.4330
ITERATION:	10	DEVIANCE:	7.4330
ITERATION:	11	DEVIANCE:	7.4330
ITERATION:	12	DEVIANCE:	7.4330
ITERATION:	13	DEVIANCE:	7.4330

ITERATIONS: 13				
REC	OBS	EXP	FREEMAN-TUKEY	PEARSON
1	2.0	1.6	.4403	.1113
2	12.0	17.5	-1.3592	1.7346
3	28.0	36.0	-1.3638	1.7751
4	28.0	35.0	-1.1909	1.3856
5	31.0	28.0	.5896	.3156
6	32.0	35.4	-.5343	.3202
7	104.0	96.5	.7722	.5822
8	206.0	197.3	.6327	.3874
9	186.0	178.7	.5558	.2952
10	102.0	105.1	-.2763	.0895

B	COEFFICIENT	STD. ERROR	WALD	SCORE
---	-------------	------------	------	-------

1	.0841	.0661	1.2724	.00013366
2	1.6407	.2179	7.5310	-.00000497
3	6.3035	.4565	13.8080	-.00000113
4	13.5241	.9642	14.0270	-.00000030
5	19.1696	1.7045	11.2466	-.00000012
6	.5907	.1255	4.7055	-.00000556

POWER: 1.00

GOODNESS-OF-FIT TESTS

STATISTIC	ESTIMATE	D.F.	PROB.
CHI-SQUARE	6.997	4	.1361
DEVIANCE	7.433	4	.1147

VARIANCE-COVARIANCE MATRIX OF COEFFICIENTS.

	B1	B2	B3	B4	B5	B6
B1	.004					
B2	.003	.047				
B3	.003	.010	.208			
B4	.003	.010	.011	.930		
B5	.003	.009	.010	.010	2.905	
B6	-.004	-.012	-.013	-.013	-.012	.016

CORRELATION MATRIX OF COEFFICIENTS.

	B1	B2	B3	B4	B5	B6
B1	1.000					
B2	.211	1.000				
B3	.111	.097	1.000			
B4	.053	.046	.024	1.000		
B5	.028	.025	.013	.006	1.000	
B6	-.490	-.431	-.226	-.107	-.057	1.000

TRACE OF THE HAT MATRIX AND ADJUSTED RESIDUALS.

RECORD	LEVERAGE	ADJ. RESID.
1	.9763	2.1668
2	.3088	-1.5841
3	.1888	-1.4792
4	.1777	-1.2981
5	.2216	.6367
6	.9318	-2.1671
7	.7680	1.5841
8	.8229	1.4792

9	.8248	1.2981
10	.7794	-.6367

DELETION RESIDUALS OR APPROXIMATE CHANGE IN COEFFICIENTS AFTER
DELETION OF EACH RECORD.

RECORD	B1	B2	B3	B4	B5	B6
1	-.9193	-.6402	-.7019	-.7049	-.6663	.8551
2	.0148	.2307	.0471	.0473	.0447	-.0573
3	.0052	.0151	.3258	.0167	.0158	-.0202
4	.0021	.0061	.0067	.5818	.0063	-.0081
5	-.0006	-.0018	-.0020	-.0020	-.5790	.0024
6	.0223	-.6403	-.7020	-.7050	-.6663	.8552
7	.0148	-.5164	.0471	.0473	.0447	-.0573
8	.0052	.0151	-1.3999	.0167	.0158	-.0202
9	.0021	.0061	.0067	-2.6924	.0063	-.0081
10	-.0006	-.0018	-.0020	-.0020	2.0342	.0024

On the first page of the output, we see that the model is indeed additive, we also see the input data and the value of the deviance goodness of fit at each iteration. The number of observed and expected deaths are given and the Freeman-Tukey and Pearson residuals as well. At convergence, the coefficient for smoking (No. 6) is 0.5907. Thus, smoking adds 0.59 deaths per 1,000 person-years (or 59/100,000) of follow-up at each age. However, since the χ^2 and D goodness of fit statistics are greater than 4 degrees of freedom, we conclude that the model does not fit the data well. Table 6 shows the comparison between the modeled baseline rates and modeled baseline rates to which 0.59/1,000 was added.

Table 6: Interpretation of absolute risks from the additive model.

Age group, i	Baseline rate, $\lambda_{i0}(\mathbf{x}'_{i0}\beta)$ (β_j) per 1,000 T	Smoker rate, $\lambda_{i1}(\mathbf{x}'_{i0}\beta)$ ($\lambda_{i0}(\mathbf{x}'_{i0}\beta) + \beta_{\text{smoke}}$) per 1,000 T
35-44	0.0841	0.6748
45-54	1.6407	2.2314
55-64	6.3035	6.8942
65-74	13.5241	14.115
75-84	19.1696	19.7603

Example 2: Multiplicative (log-linear) model using data in Table 1 adapted from [21]. The input data file for this run is the same as that in Example 1.

The screen queries to be entered by the user are as follows:

ENTER THE NAME OF INPUT FILE:
EXAMP1-3.DAT

ENTER THE NAME OF INPUT FILE:
EXAMPLE2.OUT

ENTER 1-ADD 2-MULT 3-POWER 4-NON-LINEAR
2

ENTER NUMBER OF PREDICTOR (INPUT) VARIABLES
6

NUMBER OF PARAMETERS TO BE FITTED?
6

EXPONENTIATE COEFFICIENTS? 0-NO 1-YES
1

PRINT INFORMATION AND CORR MATRICES? 0-NO 1-YES
0

PRINT INPUT DATA? 0-NO 1-YES
0

PRINT FITTED CELLS? 0-NO 1-YES
0

PERFORM REGRESSION DIAGNOSTICS? 0-NO 1-YES
0

The output file contains:

TITLE: PIRLS

MODEL SELECTED: MULTIPLICATIVE

INPUT DATA FORMAT: (F4.0,F8.3,6F3.0)

ITERATION:	1	DEVIANCE:	53.9002
ITERATION:	2	DEVIANCE:	17.9280
ITERATION:	3	DEVIANCE:	12.1932
ITERATION:	4	DEVIANCE:	12.1324
ITERATION:	5	DEVIANCE:	12.1324

B	COEFFICIENT	STD. ERROR	WALD	SCORE
1	-1.0116	.1918	-5.2751	.00000207
2	.4724	.1304	3.6236	.00000969
3	1.6159	.1147	14.0944	-.00002298
4	2.3389	.1162	20.1343	.00001905
5	2.6885	.1250	21.5110	.00001919

6 .3545 .1074 3.3019 .00002873

POWER: .00

GOODNESS-OF-FIT TESTS

STATISTIC	ESTIMATE	D.F.	PROB.
CHI-SQUARE	11.155	4	.0249
DEVIANCE	12.132	4	.0164

At convergence, the coefficient for smoking (No. 6) is 0.3545. Since we exponentiated the regression coefficients, we can interpret the RR for smoking and the 95% confidence interval as being 1.43(1.15,1.76). This implies that the risk of coronary death in smokers is 1.43 times the age-specific rates of the non-smokers. The goodness of fit statistics for this run are greater than those generated in Example 1, indicating that the multiplicative model fits worse when compared with the additive model. Table 7 shows the comparison between modeled baseline rates and the modeled baseline rates multiplied by the RR of 1.43.

Table 7: Interpretation of relative risks from the multiplicative model.

Age group, i	Baseline rate, $\lambda_{i0}(\mathbf{x}'_{i0}\beta)$ (e^{β_j}) per 1,000 T	Smoker rate, $\lambda_{i1}(\mathbf{x}'_{i0}\beta)$ ($\lambda_{i0}(\mathbf{x}'_{i0}\beta) \times e^{\beta_{\text{smoke}}}$) per 1,000 T
35-44	0.36	0.5148
45-54	1.60	2.28
55-64	5.03	7.192
65-74	10.37	14.83
75-84	14.71	21.04

Example 3: Power model using data in Table 1 adapted from [21]. The input data file for this run is the same as that in Example 1.

The screen queries to be entered by the user are as follows:

ENTER THE NAME OF INPUT FILE:
EXAMP1-3.DAT

ENTER THE NAME OF INPUT FILE:
EXAMPLE3.OUT

ENTER 1-ADD 2-MULT 3-POWER 4-NON-LINEAR
3

ENTER NUMBER OF PREDICTOR (INPUT) VARIABLES
6

NUMBER OF PARAMETERS TO BE FITTED?
6

EXPONENTIATE COEFFICIENTS? 0-NO 1-YES


```

0

PRINT INFORMATION AND CORR MATRICES?  0-NO    1-YES
0

PRINT INPUT DATA?  0-NO    1-YES
0

PRINT FITTED CELLS?  0-NO    1-YES
0

PERFORM REGRESSION DIAGNOSTICS?  0-NO    1-YES
0

```

This run fitted ten models with ρ varying from one to zero, i.e., going from a purely additive link, $\rho = 1$, to a purely multiplicative link, $\rho = 0$. The model equation is $\lambda_i(\mathbf{x}'_i\boldsymbol{\beta})^\rho$ where ρ is the power function. By inspecting the fits of the models as a function of ρ , one can see that for these example data, the lowest deviance is obtained with ρ between 0.5 to 0.6. A single run can now be made to fit the model with $\rho = 0.55$ by following the directions given in Example 1, but by specifying 0.55 for ρ instead of one. Table 8 lists the modeled baseline rates, $\beta_j^{1/0.55}$, and modeled baseline rates to which $\beta_{smoke}^{1/0.55}$ was added.

Table 8: Interpretation of relative risks from the power model.

Age group, i	Baseline rate, $\lambda_{i0}(\mathbf{x}'_{i0}\boldsymbol{\beta})$ ($\beta_j^{1/0.55}$) per 1,000 T	Smoker rate, $\lambda_{i1}(\mathbf{x}'_{i0}\boldsymbol{\beta})$ ($\lambda_{i0}(\mathbf{x}'_{i0}\boldsymbol{\beta}) + \beta_{smoke}^{1/0.55}$) per 1,000 T
35-44	0.096	0.3742
45-54	1.218	1.494
55-64	5.123	5.400
65-74	11.649	11.925
75-84	17.081	17.357

Example 4: Non-linear model using cell concentrations, and radiation doses to estimate number of splenic stem cell colonies formed after transplantation into laboratory mice. The input data (see Table 4) for this run is provided in distribution file EXAMPLE4.DAT and is listed below:

```

Till and McCulloch (1961)
(F4.0,F4.0,F7.2,F6.2)
60.  6.  1.25  .00
66.  7.  1.75  .96
46.  4.  3.00  1.92
82.  9.  7.20  2.88
105. 11. 24.00  4.32
123. 15. 75.00  5.76
12.  4. 120.00  6.72

```

The screen queries to be entered by the user are as follows:

```

ENTER THE NAME OF INPUT FILE:
EXAMPLE4.DAT

```

```

ENTER THE NAME OF INPUT FILE:

```

EXAMPLE4.OUT

ENTER 1-ADD 2-MULT 3-POWER 4-NON-LINEAR

4

ENTER NUMBER OF PREDICTOR (INPUT) VARIABLES

2

NUMBER OF PARAMETERS TO BE FITTED?

3

INPUT OR ESTIMATE STARTING VALUES: 0-INPUT 1-EST

0

INPUT STARTING VALUE FOR COEFFICIENT: 1

7.6364

INPUT STARTING VALUE FOR COEFFICIENT: 2

0.9341

INPUT STARTING VALUE FOR COEFFICIENT: 3

2.8924

PRINT INFORMATION AND CORR MATRICES? 0-NO 1-YES

1

PRINT INPUT DATA? 0-NO 1-YES

1

PRINT FITTED CELLS? 0-NO 1-YES

1

PERFORM REGRESSION DIAGNOSTICS? 0-NO 1-YES

1

Results in the output file EXAMPLE4.OUT (filename specified above) indicate that 2 iterations were required before convergence was reached. Although the deviance and chi-square GOF values were greater than the 4 d.f., the Wald statistics for the parameters were statistically significant. The scores for all coefficients were near zero. The regression parameters, standard errors (with the exception of $s.e.(\beta_1)$), correlation matrix, chi-square and deviance GOFs were identical to results in Frome [3].

TITLE: Till and McCulloch (1961)

MODEL SELECTED: NON-LINEAR

INPUT DATA FORMAT: (F4.0,F4.0,F7.2,F6.2)

60. 6. 1.25 .00

66. 7. 1.75 .96

46. 4. 3.00 1.92

82.	9.	7.20	2.88
105.	11.	24.00	4.32
123.	15.	75.00	5.76
12.	4.	120.00	6.72

STARTING VALUES FOR PARAMETERS

PARAMETER STARTING VALUE

1	7.6364000
2	.9341000
3	2.8924000

ITERATION:	1	DEVIANCE:	8.0174
ITERATION:	2	DEVIANCE:	8.0174

ITERATIONS:		2			
REC	OBS	EXP	FREEMAN-TUKEY	PEARSON	
1	60.0	57.3	.3874	.1298	
2	66.0	73.0	-.8079	.6713	
3	46.0	37.5	1.3496	1.9260	
4	82.0	91.0	-.9414	.8942	
5	105.0	101.4	.3807	.1296	
6	123.0	113.9	.8536	.7194	
7	12.0	19.9	-1.9041	3.1247	

B	COEFFICIENT	STD. ERROR	WALD	SCORE
1	7.6364	.1186	64.3725	.00030114
2	.9341	.0399	23.4229	-.00096028
3	2.8924	.7476	3.8688	.00006890

POWER: .00

GOODNESS-OF-FIT TESTS

STATISTIC	ESTIMATE	D.F.	PROB.
CHI-SQUARE	7.595	4	.1076
DEVIANCE	8.017	4	.0909

VARIANCE-COVARIANCE MATRIX OF COEFFICIENTS.

	B1	B2	B3
B1	.014		
B2	-.002	.002	

B3	-.066	.025	.559
----	-------	------	------

CORRELATION MATRIX OF COEFFICIENTS.

	B1	B2	B3
B1	1.000		
B2	-.343	1.000	
B3	-.741	.853	1.000

TRACE OF THE HAT MATRIX AND ADJUSTED RESIDUALS.

RECORD	LEVERAGE	ADJ. RESID.
1	.8060	.8181
2	.3663	-1.0293
3	.2251	1.5766
4	.4551	-1.2811
5	.2954	.4289
6	.6358	1.4055
7	.2162	-1.9966

DELETION RESIDUALS OR APPROXIMATE CHANGE IN COEFFICIENTS AFTER DELETION OF EACH RECORD.

RECORD	B1	B2	B3
1	-.1978	.0228	.9235
2	.0581	.0141	.0336
3	.0065	-.0276	-.4434
4	-.0339	.0336	.6643
5	.0060	-.0010	-.0749
6	-.0201	.0499	.4879
7	.0231	-.0354	-.4386